



Full wwPDB X-ray Structure Validation Report ⓘ

Sep 24, 2020 – 02:19 PM EDT

PDB ID : 6V75
Title : Crystal Structure of Human PKM2 in Complex with L-aspartate
Authors : Nandi, S.; Dey, M.
Deposited on : 2019-12-07
Resolution : 2.85 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.14.6
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.14.6

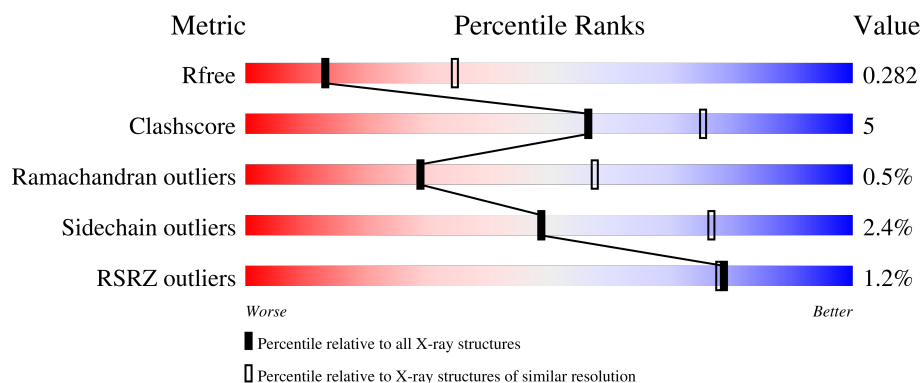
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.85 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	3168 (2.90-2.82)
Clashscore	141614	3438 (2.90-2.82)
Ramachandran outliers	138981	3348 (2.90-2.82)
Sidechain outliers	138945	3351 (2.90-2.82)
RSRZ outliers	127900	3103 (2.90-2.82)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	550	<div> <div>%</div> <div>82% 10% 7%</div> </div>
1	B	550	<div> <div>2%</div> <div>80% 13% 7%</div> </div>
1	C	550	<div> <div>%</div> <div>84% 8% 6%</div> </div>
1	D	550	<div> <div>%</div> <div>69% 9% 22%</div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard

residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
6	MG	A	607	-	-	-	X

2 Entry composition

There are 10 unique types of molecules in this entry. The entry contains 13674 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called Pyruvate kinase PKM.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	C	517	Total	C	N	O	S	0	0	0
			3550	2223	632	671	24			
1	A	509	Total	C	N	O	S	0	0	0
			3450	2152	608	667	23			
1	B	512	Total	C	N	O	S	0	0	0
			3491	2166	632	671	22			
1	D	430	Total	C	N	O	S	0	0	0
			2917	1809	525	563	20			

There are 76 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	-18	MET	-	initiating methionine	UNP P14618
C	-17	GLY	-	expression tag	UNP P14618
C	-16	SER	-	expression tag	UNP P14618
C	-15	SER	-	expression tag	UNP P14618
C	-14	HIS	-	expression tag	UNP P14618
C	-13	HIS	-	expression tag	UNP P14618
C	-12	HIS	-	expression tag	UNP P14618
C	-11	HIS	-	expression tag	UNP P14618
C	-10	HIS	-	expression tag	UNP P14618
C	-9	HIS	-	expression tag	UNP P14618
C	-8	SER	-	expression tag	UNP P14618
C	-7	SER	-	expression tag	UNP P14618
C	-6	GLY	-	expression tag	UNP P14618
C	-5	LEU	-	expression tag	UNP P14618
C	-4	VAL	-	expression tag	UNP P14618
C	-3	PRO	-	expression tag	UNP P14618
C	-2	ARG	-	expression tag	UNP P14618
C	-1	GLY	-	expression tag	UNP P14618
C	0	SER	-	expression tag	UNP P14618
A	-18	MET	-	initiating methionine	UNP P14618
A	-17	GLY	-	expression tag	UNP P14618

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Chain	Residue	Modelled	Actual	Comment	Reference
A	-16	SER	-	expression tag	UNP P14618
A	-15	SER	-	expression tag	UNP P14618
A	-14	HIS	-	expression tag	UNP P14618
A	-13	HIS	-	expression tag	UNP P14618
A	-12	HIS	-	expression tag	UNP P14618
A	-11	HIS	-	expression tag	UNP P14618
A	-10	HIS	-	expression tag	UNP P14618
A	-9	HIS	-	expression tag	UNP P14618
A	-8	SER	-	expression tag	UNP P14618
A	-7	SER	-	expression tag	UNP P14618
A	-6	GLY	-	expression tag	UNP P14618
A	-5	LEU	-	expression tag	UNP P14618
A	-4	VAL	-	expression tag	UNP P14618
A	-3	PRO	-	expression tag	UNP P14618
A	-2	ARG	-	expression tag	UNP P14618
A	-1	GLY	-	expression tag	UNP P14618
A	0	SER	-	expression tag	UNP P14618
B	-18	MET	-	initiating methionine	UNP P14618
B	-17	GLY	-	expression tag	UNP P14618
B	-16	SER	-	expression tag	UNP P14618
B	-15	SER	-	expression tag	UNP P14618
B	-14	HIS	-	expression tag	UNP P14618
B	-13	HIS	-	expression tag	UNP P14618
B	-12	HIS	-	expression tag	UNP P14618
B	-11	HIS	-	expression tag	UNP P14618
B	-10	HIS	-	expression tag	UNP P14618
B	-9	HIS	-	expression tag	UNP P14618
B	-8	SER	-	expression tag	UNP P14618
B	-7	SER	-	expression tag	UNP P14618
B	-6	GLY	-	expression tag	UNP P14618
B	-5	LEU	-	expression tag	UNP P14618
B	-4	VAL	-	expression tag	UNP P14618
B	-3	PRO	-	expression tag	UNP P14618
B	-2	ARG	-	expression tag	UNP P14618
B	-1	GLY	-	expression tag	UNP P14618
B	0	SER	-	expression tag	UNP P14618
D	-18	MET	-	initiating methionine	UNP P14618
D	-17	GLY	-	expression tag	UNP P14618
D	-16	SER	-	expression tag	UNP P14618
D	-15	SER	-	expression tag	UNP P14618
D	-14	HIS	-	expression tag	UNP P14618
D	-13	HIS	-	expression tag	UNP P14618

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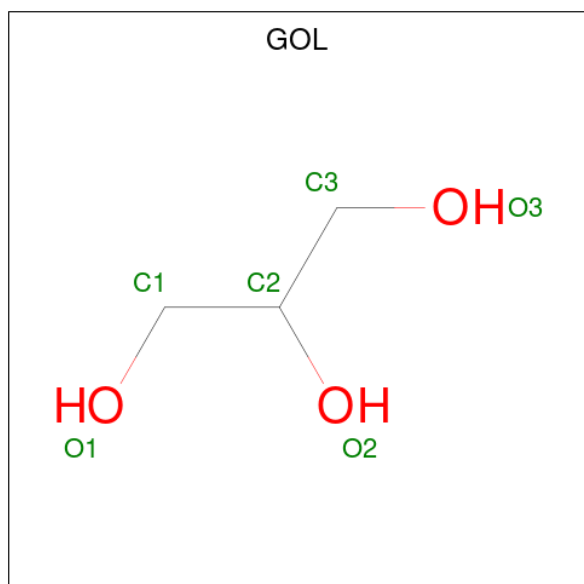
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Chain	Residue	Modelled	Actual	Comment	Reference
D	-12	HIS	-	expression tag	UNP P14618
D	-11	HIS	-	expression tag	UNP P14618
D	-10	HIS	-	expression tag	UNP P14618
D	-9	HIS	-	expression tag	UNP P14618
D	-8	SER	-	expression tag	UNP P14618
D	-7	SER	-	expression tag	UNP P14618
D	-6	GLY	-	expression tag	UNP P14618
D	-5	LEU	-	expression tag	UNP P14618
D	-4	VAL	-	expression tag	UNP P14618
D	-3	PRO	-	expression tag	UNP P14618
D	-2	ARG	-	expression tag	UNP P14618
D	-1	GLY	-	expression tag	UNP P14618
D	0	SER	-	expression tag	UNP P14618

- Molecule 2 is POTASSIUM ION (three-letter code: K) (formula: K).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	B	1	Total K 1 1	0	0
2	A	1	Total K 1 1	0	0
2	D	2	Total K 2 2	0	0
2	C	2	Total K 2 2	0	0

- Molecule 3 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).

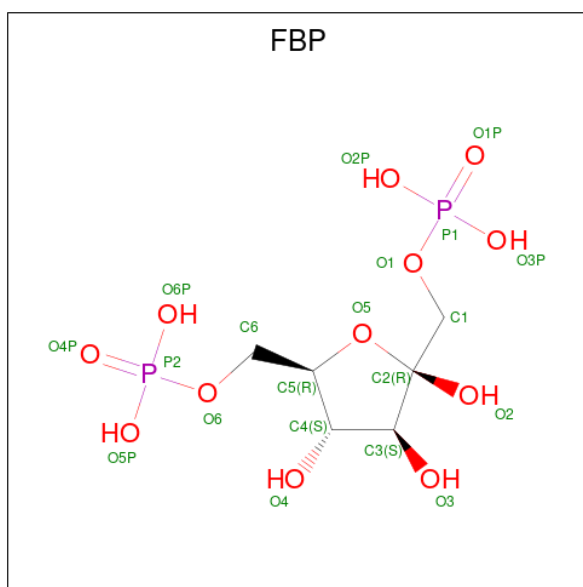


Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	C	1	Total C O 6 3 3	0	0
3	C	1	Total C O 6 3 3	0	0
3	C	1	Total C O 6 3 3	0	0
3	A	1	Total C O 6 3 3	0	0
3	A	1	Total C O 6 3 3	0	0
3	B	1	Total C O 6 3 3	0	0
3	B	1	Total C O 6 3 3	0	0
3	B	1	Total C O 6 3 3	0	0
3	B	1	Total C O 6 3 3	0	0
3	B	1	Total C O 6 3 3	0	0
3	D	1	Total C O 6 3 3	0	0
3	D	1	Total C O 6 3 3	0	0

- Molecule 4 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	C	1	Total Cl 1 1	0	0

- Molecule 5 is 1,6-di-O-phosphono-beta-D-fructofuranose (three-letter code: FBP) (formula: C₆H₁₄O₁₂P₂).

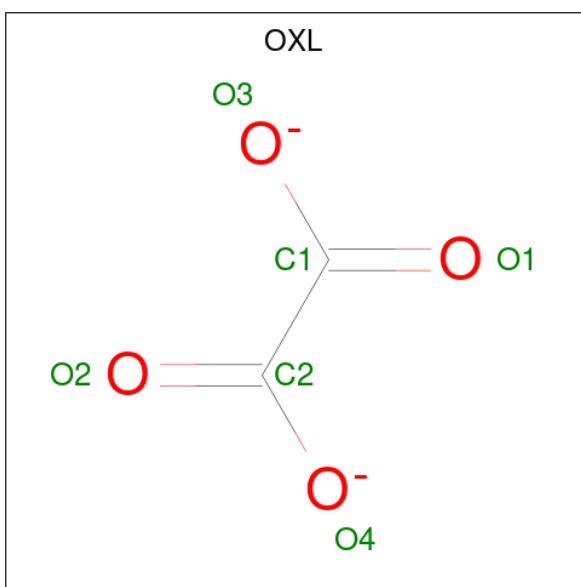


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	C	1	Total	C	O	P	0	0
			20	6	12	2		
5	A	1	Total	C	O	P	0	0
			20	6	12	2		
5	B	1	Total	C	O	P	0	0
			20	6	12	2		
5	D	1	Total	C	O	P	0	0
			20	6	12	2		

- Molecule 6 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

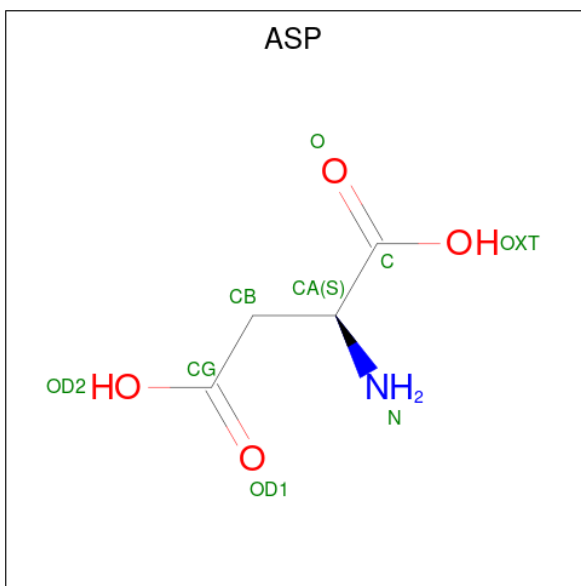
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	B	1	Total	Mg	0	0
			1	1		
6	A	1	Total	Mg	0	0
			1	1		
6	D	1	Total	Mg	0	0
			1	1		
6	C	1	Total	Mg	0	0
			1	1		

- Molecule 7 is OXALATE ION (three-letter code: OXL) (formula: C₂O₄).



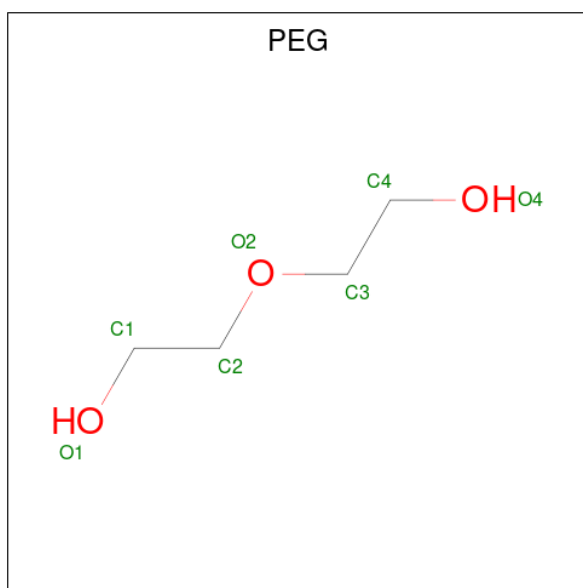
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	C	1	Total	C	O	0	0
			6	2	4		

- Molecule 8 is ASPARTIC ACID (three-letter code: ASP) (formula: $C_4H_7NO_4$) (labeled as "Ligand of Interest" by author).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
8	C	1	Total	C	N	O	0	0
			9	4	1	4		
8	B	1	Total	C	N	O	0	0
			9	4	1	4		

- Molecule 9 is DI(HYDROXYETHYL)ETHER (three-letter code: PEG) (formula: $C_4H_{10}O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
9	A	1	Total	C	O	0	0
			7	4	3		
9	A	1	Total	C	O	0	0
			7	4	3		
9	B	1	Total	C	O	0	0
			7	4	3		

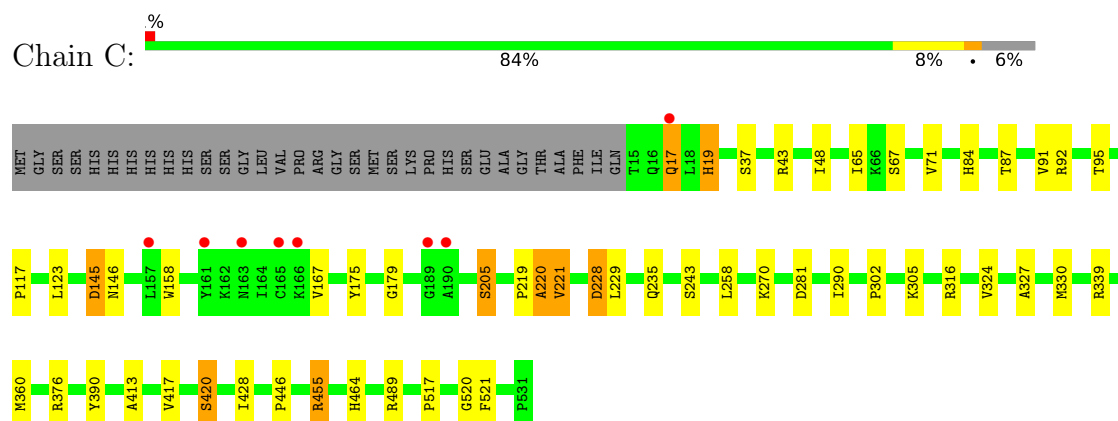
- Molecule 10 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
10	C	15	Total	O	0	0
			15	15		
10	A	24	Total	O	0	0
			24	24		
10	B	7	Total	O	0	0
			7	7		
10	D	12	Total	O	0	0
			12	12		

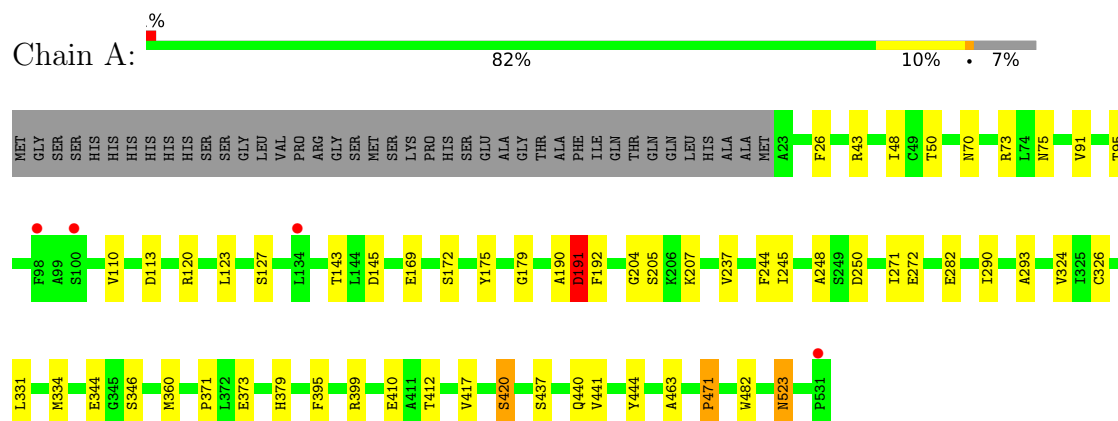
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

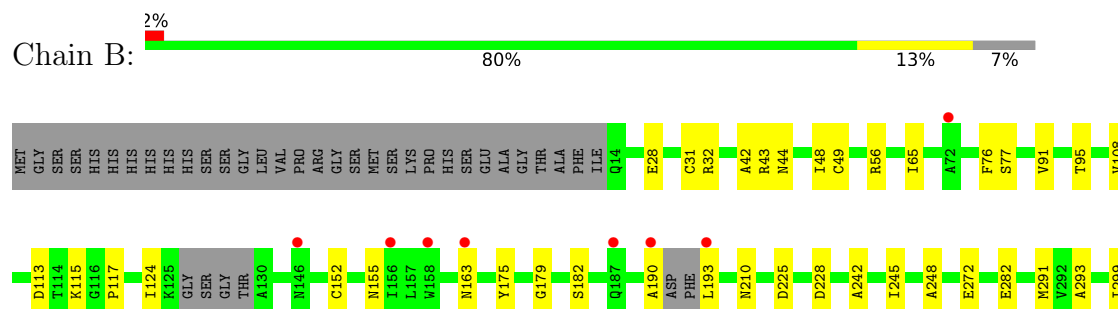
• Molecule 1: Pyruvate kinase PKM

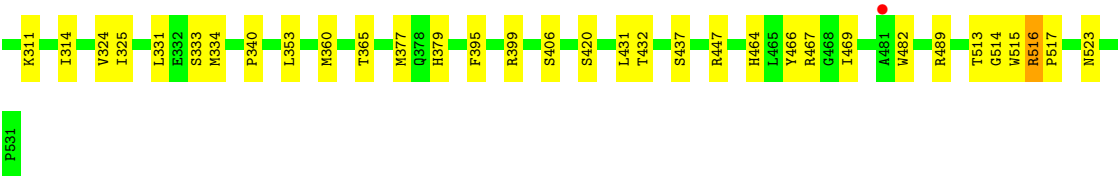


• Molecule 1: Pyruvate kinase PKM

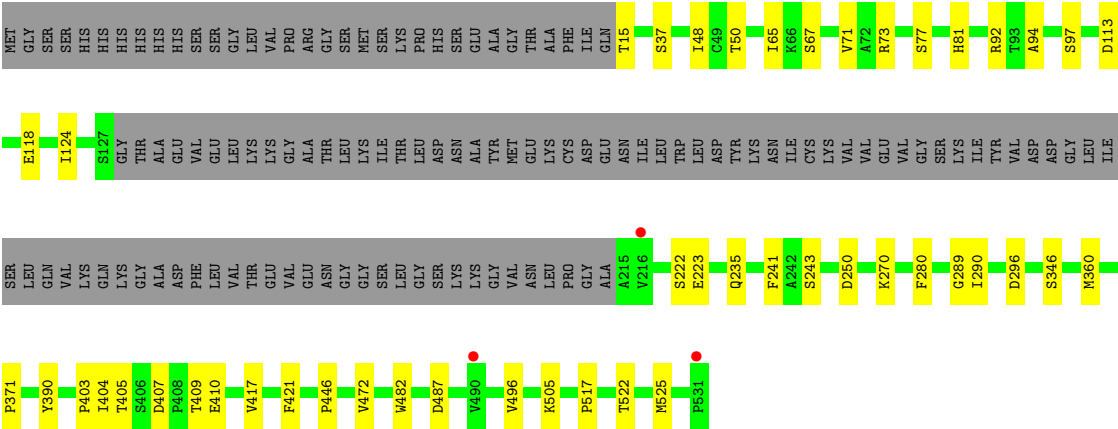


• Molecule 1: Pyruvate kinase PKM





• Molecule 1: Pyruvate kinase PKM



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	82.52Å 157.96Å 93.03Å 90.00° 101.08° 90.00°	Depositor
Resolution (Å)	55.52 – 2.85 55.52 – 2.85	Depositor EDS
% Data completeness (in resolution range)	99.3 (55.52-2.85) 99.4 (55.52-2.85)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.97 (at 2.86Å)	Xtriage
Refinement program	PHENIX 1.17.1_3660	Depositor
R, R_{free}	0.246 , 0.284 0.245 , 0.282	Depositor DCC
R_{free} test set	2637 reflections (4.86%)	wwPDB-VP
Wilson B-factor (Å ²)	69.0	Xtriage
Anisotropy	0.076	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 82.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	13674	wwPDB-VP
Average B, all atoms (Å ²)	67.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.60% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: GOL, MG, FBP, CL, K, OXL, PEG

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.25	0/3509	0.49	3/4785 (0.1%)
1	B	0.28	0/3548	0.47	2/4832 (0.0%)
1	C	0.27	0/3611	0.47	5/4923 (0.1%)
1	D	0.24	0/2967	0.45	2/4047 (0.0%)
All	All	0.26	0/13635	0.47	12/18587 (0.1%)

There are no bond length outliers.

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	155	ASN	CB-CA-C	8.47	127.34	110.40
1	A	523	ASN	CB-CA-C	8.34	127.08	110.40
1	C	327	ALA	N-CA-C	6.92	129.68	111.00
1	A	191	ASP	N-CA-C	-6.40	93.72	111.00
1	D	505	LYS	CB-CA-C	5.88	122.16	110.40
1	C	327	ALA	CB-CA-C	-5.76	101.46	110.10
1	C	145	ASP	CB-CA-C	-5.57	99.26	110.40
1	C	220	ALA	CB-CA-C	-5.46	101.91	110.10
1	B	523	ASN	CB-CA-C	5.30	121.00	110.40
1	A	244	PHE	N-CA-C	5.21	125.06	111.00
1	C	146	ASN	CB-CA-C	5.12	120.64	110.40
1	D	505	LYS	N-CA-C	-5.04	97.38	111.00

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3450	0	2990	31	0
1	B	3491	0	3021	43	0
1	C	3550	0	3161	32	0
1	D	2917	0	2524	22	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	2	0	0	0	0
2	D	2	0	0	0	0
3	A	12	0	16	0	0
3	B	30	0	40	4	0
3	C	18	0	24	0	0
3	D	12	0	16	1	0
4	C	1	0	0	0	0
5	A	20	0	10	0	0
5	B	20	0	10	3	0
5	C	20	0	10	4	0
5	D	20	0	10	0	0
6	A	1	0	0	0	0
6	B	1	0	0	0	0
6	C	1	0	0	0	0
6	D	1	0	0	0	0
7	C	6	0	0	0	0
8	B	9	0	3	3	0
8	C	9	0	3	1	0
9	A	14	0	20	2	0
9	B	7	0	10	0	0
10	A	24	0	0	1	0
10	B	7	0	0	1	0
10	C	15	0	0	0	0
10	D	12	0	0	0	0
All	All	13674	0	11868	127	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 5.

All (127) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:489:ARG:NH2	5:C:607:FBP:O2P	1.70	1.24
1:C:19:HIS:ND1	1:C:19:HIS:O	2.11	0.83
1:B:469:ILE:O	8:B:610:ASP:HB2	1.81	0.80
1:B:95:THR:HG21	1:B:108:VAL:HG21	1.62	0.80
1:A:127:SER:OG	10:A:701:HOH:O	2.04	0.75
1:A:123:LEU:HA	1:A:205:SER:HB3	1.70	0.74
1:A:412:THR:HG21	1:A:523:ASN:O	1.87	0.74
1:C:175:TYR:HB3	1:C:179:GLY:HA2	1.69	0.73
1:A:412:THR:CG2	1:A:523:ASN:O	2.41	0.69
1:D:407:ASP:HB3	1:D:410:GLU:HB2	1.75	0.69
1:C:417:VAL:HG13	1:C:446:PRO:HB3	1.76	0.67
1:A:43:ARG:NH2	1:A:70:ASN:OD1	2.28	0.67
1:C:145:ASP:O	1:C:158:TRP:CD1	2.49	0.66
1:D:472:VAL:HG11	1:D:496:VAL:HG11	1.77	0.66
1:C:290:ILE:HB	1:C:324:VAL:HG12	1.77	0.66
1:A:143:THR:HG22	1:A:145:ASP:H	1.60	0.65
1:B:464:HIS:HD1	8:B:610:ASP:N	1.96	0.63
1:D:50:THR:OG1	1:D:73:ARG:NH1	2.31	0.63
1:B:190:ALA:C	1:B:193:LEU:N	2.53	0.62
1:B:311:LYS:NZ	3:B:603:GOL:O3	2.32	0.62
1:A:48:ILE:HB	1:A:360:MET:HG3	1.82	0.62
1:C:517:PRO:HA	5:C:607:FBP:H11	1.80	0.61
1:B:489:ARG:NH2	5:B:607:FBP:O1P	2.33	0.61
1:A:271:ILE:HD11	1:A:290:ILE:HD12	1.83	0.60
1:C:92:ARG:NH2	1:C:235:GLN:O	2.34	0.60
1:D:48:ILE:HB	1:D:360:MET:HG3	1.84	0.60
1:A:272:GLU:HG2	1:A:293:ALA:HB3	1.84	0.60
1:C:464:HIS:HD1	8:C:610:ASP:N	2.01	0.59
1:D:92:ARG:NH1	1:D:235:GLN:O	2.34	0.58
1:D:94:ALA:O	1:D:97:SER:OG	2.20	0.58
9:A:605:PEG:H11	1:B:56:ARG:HH11	1.69	0.58
1:C:520:GLY:N	5:C:607:FBP:O4P	2.35	0.57
1:B:466:TYR:O	8:B:610:ASP:OD1	2.24	0.56
1:D:417:VAL:HG13	1:D:446:PRO:HB3	1.88	0.56
1:C:67:SER:OG	1:C:376:ARG:NH1	2.39	0.56
1:A:417:VAL:HA	1:A:420:SER:HB3	1.89	0.55
1:A:482:TRP:HB3	9:A:605:PEG:H21	1.90	0.54
1:A:463:ALA:HB3	1:A:471:PRO:HG3	1.89	0.54
1:B:225:ASP:HA	1:B:228:ASP:HB2	1.90	0.54
1:B:48:ILE:HB	1:B:360:MET:HG3	1.89	0.53
1:D:48:ILE:HG23	1:D:71:VAL:HG13	1.91	0.53
1:B:482:TRP:CD1	1:B:517:PRO:HG3	2.44	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:42:ALA:HB1	3:B:602:GOL:H2	1.91	0.52
1:B:466:TYR:HB2	1:B:469:ILE:HD12	1.91	0.52
1:C:521:PHE:HB2	5:C:607:FBP:O4	2.09	0.52
1:B:49:CYS:HB3	1:B:365:THR:HG21	1.91	0.51
1:A:91:VAL:O	1:A:95:THR:HG23	2.11	0.51
1:C:281:ASP:OD1	1:C:316:ARG:NH1	2.44	0.51
1:B:210:ASN:HD22	1:B:299:ILE:HG21	1.77	0.50
1:B:77:SER:HA	1:B:115:LYS:HD3	1.94	0.50
1:A:331:LEU:HD23	1:A:344:GLU:HB3	1.92	0.50
1:B:272:GLU:HG2	1:B:293:ALA:HB3	1.93	0.50
1:A:248:ALA:HB2	1:A:282:GLU:HG2	1.94	0.50
1:B:28:GLU:HG3	1:B:32:ARG:HD2	1.94	0.49
1:B:432:THR:HA	5:B:607:FBP:H61	1.93	0.49
1:D:113:ASP:HA	1:D:241:PHE:HB2	1.93	0.49
1:B:248:ALA:HB2	1:B:282:GLU:HG2	1.94	0.49
1:C:91:VAL:O	1:C:95:THR:HG23	2.12	0.49
1:D:77:SER:OG	1:D:118:GLU:OE2	2.30	0.49
1:B:515:TRP:CE3	1:B:516:ARG:HB3	2.48	0.48
1:C:48:ILE:HG12	1:C:71:VAL:HB	1.94	0.48
1:A:204:GLY:HA3	1:A:207:LYS:HE2	1.95	0.48
1:D:222:SER:OG	1:D:223:GLU:N	2.46	0.48
1:A:169:GLU:O	1:A:172:SER:OG	2.28	0.48
1:D:409:THR:HG23	1:D:522:THR:HB	1.95	0.48
1:A:245:ILE:HG23	1:A:250:ASP:HB2	1.96	0.48
1:C:145:ASP:O	1:C:158:TRP:NE1	2.46	0.47
1:B:331:LEU:HB3	1:B:334:MET:HG3	1.96	0.47
1:B:447:ARG:NH1	10:B:701:HOH:O	2.47	0.47
1:D:15:THR:O	1:D:37:SER:OG	2.31	0.47
1:C:420:SER:HB2	1:C:428:ILE:HD11	1.96	0.47
1:A:395:PHE:HE1	1:A:444:TYR:HB3	1.80	0.47
1:A:175:TYR:HB3	1:A:179:GLY:HA2	1.95	0.47
1:B:340:PRO:HG3	1:B:377:MET:HG2	1.97	0.46
1:D:65:ILE:C	1:D:67:SER:H	2.18	0.46
1:B:43:ARG:HB2	3:B:602:GOL:H31	1.97	0.46
1:D:404:ILE:O	1:D:405:THR:C	2.53	0.46
1:A:43:ARG:HD3	1:A:379:HIS:ND1	2.30	0.46
1:B:91:VAL:O	1:B:95:THR:HG23	2.15	0.46
1:D:15:THR:O	1:D:15:THR:OG1	2.27	0.46
1:C:243:SER:HA	1:C:270:LYS:HB2	1.97	0.45
1:C:290:ILE:O	1:C:324:VAL:HA	2.16	0.45
1:A:75:ASN:HA	1:A:113:ASP:HB3	1.97	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:437:SER:OG	5:B:607:FBP:O4P	2.21	0.45
1:D:241:PHE:HB3	1:D:270:LYS:HD3	1.98	0.45
1:A:395:PHE:CZ	1:A:399:ARG:HD3	2.52	0.45
1:B:431:LEU:HD21	1:B:489:ARG:HB2	1.98	0.45
1:C:117:PRO:HD3	1:C:220:ALA:HA	1.99	0.45
1:B:117:PRO:O	3:B:604:GOL:O1	2.24	0.44
1:B:190:ALA:H	1:B:193:LEU:N	2.16	0.44
1:D:280:PHE:HE1	1:D:290:ILE:HG21	1.82	0.44
1:B:175:TYR:HB3	1:B:179:GLY:HA2	1.98	0.44
1:B:65:ILE:HD11	1:B:95:THR:HG22	1.98	0.44
1:C:65:ILE:HD11	1:C:95:THR:HG22	2.00	0.44
1:B:43:ARG:HD3	1:B:379:HIS:ND1	2.32	0.44
1:D:482:TRP:CD2	1:D:517:PRO:HG3	2.53	0.43
1:B:124:ILE:HA	1:B:152:CYS:H	1.83	0.43
1:A:331:LEU:HB3	1:A:334:MET:HG3	2.00	0.43
1:C:316:ARG:HG2	1:B:31:CYS:HB3	2.00	0.43
1:C:228:ASP:N	1:C:228:ASP:OD1	2.50	0.43
1:C:302:PRO:HG2	1:C:305:LYS:HE3	1.99	0.43
1:C:330:MET:HE3	1:C:360:MET:H	1.84	0.43
1:C:413:ALA:O	1:C:417:VAL:HG23	2.19	0.43
1:A:373:GLU:N	1:A:373:GLU:OE1	2.46	0.42
1:C:123:LEU:HA	1:C:205:SER:OG	2.20	0.42
1:D:289:GLY:O	1:D:290:ILE:HD13	2.19	0.42
1:D:421:PHE:HB3	3:D:604:GOL:H2	2.02	0.42
1:B:314:ILE:HA	1:B:324:VAL:HG21	2.02	0.42
1:C:84:HIS:HA	1:C:87:THR:HG22	2.01	0.42
1:B:44:ASN:O	1:B:467:ARG:HG3	2.18	0.42
1:A:110:VAL:HG12	1:A:237:VAL:HG23	2.01	0.42
1:B:395:PHE:CZ	1:B:399:ARG:HD3	2.55	0.42
1:B:242:ALA:HB1	1:B:245:ILE:HD11	2.02	0.41
1:A:437:SER:O	1:A:441:VAL:HG23	2.19	0.41
1:B:353:LEU:HD23	1:B:353:LEU:HA	1.87	0.41
1:B:76:PHE:HB2	1:B:113:ASP:O	2.19	0.41
1:C:455:ARG:H	1:C:455:ARG:HG3	1.60	0.41
1:A:50:THR:OG1	1:A:73:ARG:NH1	2.40	0.41
1:B:291:MET:HG3	1:B:325:ILE:HB	2.03	0.41
1:A:190:ALA:O	1:A:191:ASP:CB	2.68	0.41
1:C:17:GLN:N	1:C:37:SER:OG	2.54	0.41
1:B:513:THR:OG1	1:B:514:GLY:N	2.54	0.41
1:A:290:ILE:O	1:A:324:VAL:HA	2.21	0.41
1:C:229:LEU:HD22	1:C:258:LEU:HD21	2.03	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:243:SER:HA	1:D:270:LYS:HE2	2.03	0.40
1:C:219:PRO:C	1:C:221:VAL:H	2.23	0.40
1:A:410:GLU:HG2	1:A:440:GLN:OE1	2.21	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	507/550 (92%)	471 (93%)	33 (6%)	3 (1%)	25	53
1	B	506/550 (92%)	472 (93%)	33 (6%)	1 (0%)	47	75
1	C	515/550 (94%)	478 (93%)	34 (7%)	3 (1%)	25	53
1	D	426/550 (78%)	389 (91%)	34 (8%)	3 (1%)	22	50
All	All	1954/2200 (89%)	1810 (93%)	134 (7%)	10 (0%)	29	57

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	191	ASP
1	A	192	PHE
1	D	124	ILE
1	C	167	VAL
1	A	371	PRO
1	D	371	PRO
1	C	17	GLN
1	B	163	ASN
1	D	403	PRO
1	C	221	VAL

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	283/452 (63%)	277 (98%)	6 (2%)	53	79
1	B	282/452 (62%)	277 (98%)	5 (2%)	59	82
1	C	300/452 (66%)	292 (97%)	8 (3%)	44	74
1	D	238/452 (53%)	231 (97%)	7 (3%)	42	72
All	All	1103/1808 (61%)	1077 (98%)	26 (2%)	49	77

All (26) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	19	HIS
1	C	43	ARG
1	C	205	SER
1	C	228	ASP
1	C	339	ARG
1	C	390	TYR
1	C	420	SER
1	C	455	ARG
1	A	26	PHE
1	A	120	ARG
1	A	326	CYS
1	A	346	SER
1	A	420	SER
1	A	471	PRO
1	B	182	SER
1	B	333	SER
1	B	406	SER
1	B	420	SER
1	B	516	ARG
1	D	81	HIS
1	D	250	ASP
1	D	296	ASP
1	D	346	SER
1	D	390	TYR

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Mol	Chain	Res	Type
1	D	487	ASP
1	D	525	MET

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	210	ASN
1	B	310	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 33 ligands modelled in this entry, 11 are monoatomic - leaving 22 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
3	GOL	B	603	-	5,5,5	0.91	0	5,5,5	0.95	0
3	GOL	C	603	-	5,5,5	0.92	0	5,5,5	1.01	0
3	GOL	A	603	-	5,5,5	0.93	0	5,5,5	0.93	0
3	GOL	B	606	-	5,5,5	0.92	0	5,5,5	0.98	0
9	PEG	A	606	-	6,6,6	0.48	0	5,5,5	0.46	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	FBP	C	607	-	18,20,20	1.90	6 (33%)	23,32,32	1.67	8 (34%)
5	FBP	A	604	-	18,20,20	0.93	1 (5%)	23,32,32	0.65	0
3	GOL	B	604	-	5,5,5	0.90	0	5,5,5	1.04	0
3	GOL	C	604	-	5,5,5	0.94	0	5,5,5	1.00	0
9	PEG	A	605	-	6,6,6	0.50	0	5,5,5	0.46	0
3	GOL	A	602	-	5,5,5	0.93	0	5,5,5	0.98	0
3	GOL	D	604	-	5,5,5	0.93	0	5,5,5	0.97	0
5	FBP	D	605	-	18,20,20	0.92	1 (5%)	23,32,32	0.74	0
3	GOL	B	602	-	5,5,5	0.87	0	5,5,5	1.01	0
5	FBP	B	607	-	18,20,20	0.92	1 (5%)	23,32,32	0.85	0
9	PEG	B	608	-	6,6,6	0.48	0	5,5,5	0.39	0
7	OXL	C	609	6	0,5,5	0.00	-	0,6,6	0.00	-
3	GOL	D	603	-	5,5,5	0.88	0	5,5,5	0.99	0
8	ASP	C	610	-	2,8,8	0.38	0	1,10,10	0.41	0
3	GOL	B	605	-	5,5,5	0.92	0	5,5,5	0.99	0
3	GOL	C	605	-	5,5,5	0.91	0	5,5,5	1.01	0
8	ASP	B	610	-	2,8,8	0.31	0	1,10,10	0.38	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	GOL	B	603	-	-	2/4/4/4	-
3	GOL	C	603	-	-	0/4/4/4	-
3	GOL	A	603	-	-	2/4/4/4	-
3	GOL	B	606	-	-	0/4/4/4	-
9	PEG	A	606	-	-	0/4/4/4	-
5	FBP	C	607	-	-	5/13/32/32	0/1/1/1
5	FBP	A	604	-	-	4/13/32/32	0/1/1/1
3	GOL	B	604	-	-	1/4/4/4	-
3	GOL	C	604	-	-	0/4/4/4	-
9	PEG	A	605	-	-	1/4/4/4	-
3	GOL	A	602	-	-	0/4/4/4	-
3	GOL	D	604	-	-	2/4/4/4	-
5	FBP	D	605	-	-	7/13/32/32	0/1/1/1
3	GOL	B	602	-	-	2/4/4/4	-
5	FBP	B	607	-	-	11/13/32/32	0/1/1/1
9	PEG	B	608	-	-	2/4/4/4	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
7	OXL	C	609	6	-	0/0/4/4	-
3	GOL	D	603	-	-	2/4/4/4	-
8	ASP	C	610	-	-	1/2/8/8	-
3	GOL	B	605	-	-	0/4/4/4	-
3	GOL	C	605	-	-	4/4/4/4	-
8	ASP	B	610	-	-	1/2/8/8	-

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	C	607	FBP	P2-O6P	-3.37	1.41	1.54
5	C	607	FBP	P2-O5P	-3.10	1.42	1.54
5	C	607	FBP	P1-O3P	-2.77	1.44	1.54
5	B	607	FBP	O2-C2	2.75	1.45	1.40
5	A	604	FBP	O2-C2	2.71	1.45	1.40
5	D	605	FBP	O2-C2	2.70	1.45	1.40
5	C	607	FBP	O5-C2	-2.56	1.39	1.43
5	C	607	FBP	P1-O2P	-2.28	1.46	1.54
5	C	607	FBP	P2-O6	-2.06	1.53	1.60

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	C	607	FBP	O6P-P2-O6	-3.02	98.70	106.73
5	C	607	FBP	P2-O6-C6	2.69	125.71	118.30
5	C	607	FBP	O2P-P1-O1P	2.58	120.78	110.68
5	C	607	FBP	O3-C3-C4	-2.29	105.40	113.32
5	C	607	FBP	O2-C2-O5	-2.20	105.26	109.50
5	C	607	FBP	O5-C5-C6	2.08	114.03	109.45
5	C	607	FBP	O4-C4-C3	-2.03	106.09	112.15
5	C	607	FBP	O6P-P2-O4P	2.02	118.58	110.68

There are no chirality outliers.

All (47) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	603	GOL	O1-C1-C2-C3
3	A	603	GOL	C1-C2-C3-O3
5	C	607	FBP	O1-C1-C2-O2
5	C	607	FBP	O1-C1-C2-C3
5	C	607	FBP	O1-C1-C2-O5

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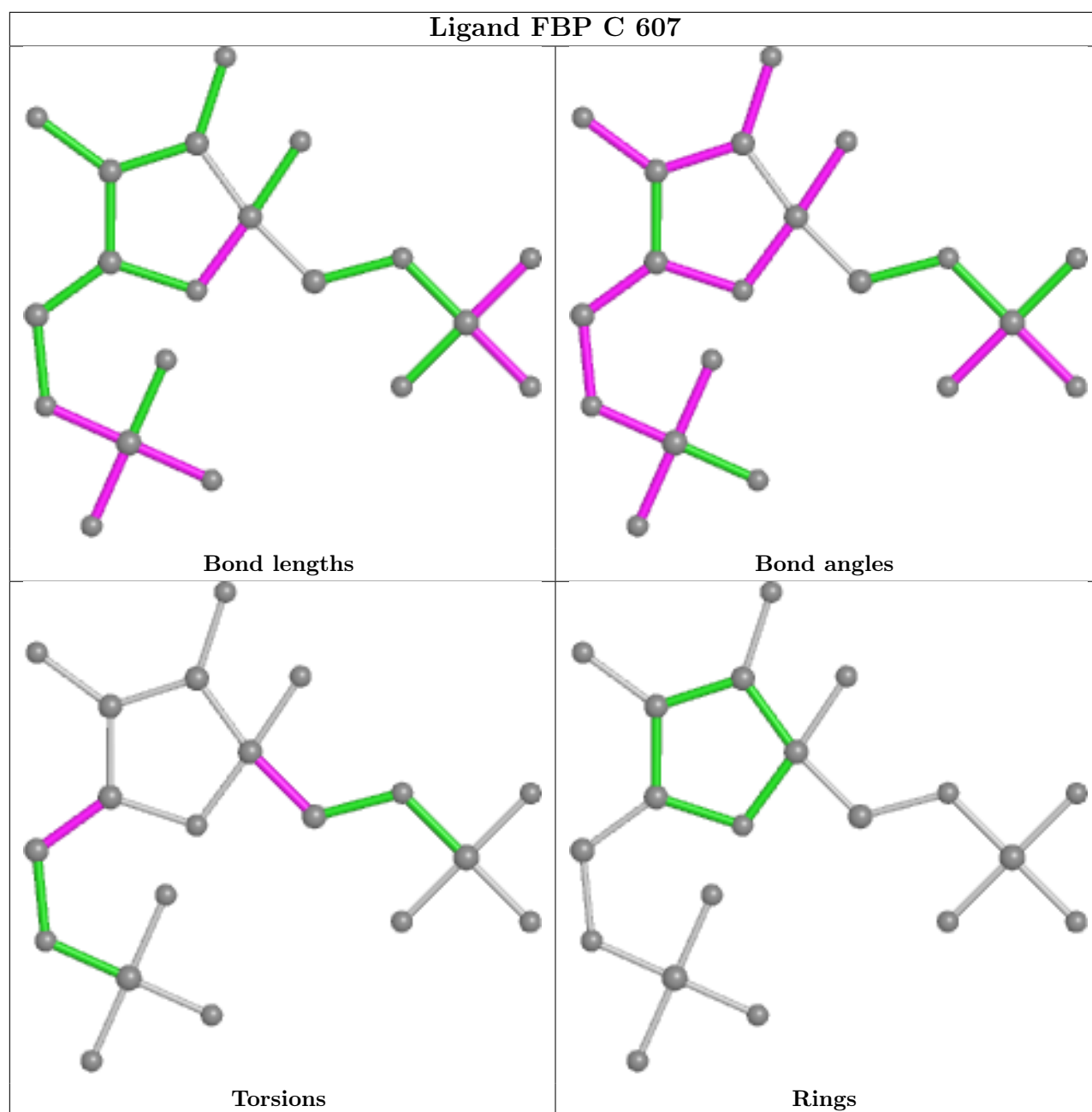
Mol	Chain	Res	Type	Atoms
5	C	607	FBP	C4-C5-C6-O6
5	A	604	FBP	O1-C1-C2-O2
5	A	604	FBP	O1-C1-C2-C3
5	A	604	FBP	O1-C1-C2-O5
3	C	605	GOL	O1-C1-C2-O2
3	C	605	GOL	O1-C1-C2-C3
3	C	605	GOL	C1-C2-C3-O3
3	D	604	GOL	O1-C1-C2-C3
5	D	605	FBP	O1-C1-C2-O2
5	D	605	FBP	O1-C1-C2-C3
5	D	605	FBP	O1-C1-C2-O5
5	D	605	FBP	C6-O6-P2-O5P
5	D	605	FBP	C6-O6-P2-O6P
3	B	602	GOL	O1-C1-C2-C3
5	B	607	FBP	C1-O1-P1-O1P
5	B	607	FBP	C1-O1-P1-O2P
5	B	607	FBP	C1-O1-P1-O3P
5	B	607	FBP	O1-C1-C2-O2
5	B	607	FBP	O1-C1-C2-C3
5	B	607	FBP	O1-C1-C2-O5
5	B	607	FBP	C4-C5-C6-O6
5	B	607	FBP	C6-O6-P2-O4P
5	B	607	FBP	C6-O6-P2-O5P
5	B	607	FBP	C6-O6-P2-O6P
3	D	603	GOL	O1-C1-C2-O2
3	D	603	GOL	O1-C1-C2-C3
8	C	610	ASP	N-CA-CB-CG
3	C	605	GOL	O2-C2-C3-O3
5	B	607	FBP	O5-C5-C6-O6
9	B	608	PEG	O2-C3-C4-O4
3	B	602	GOL	O1-C1-C2-O2
5	C	607	FBP	O5-C5-C6-O6
8	B	610	ASP	N-CA-CB-CG
3	D	604	GOL	O1-C1-C2-O2
5	D	605	FBP	C6-O6-P2-O4P
3	A	603	GOL	O2-C2-C3-O3
9	A	605	PEG	C4-C3-O2-C2
3	B	603	GOL	O1-C1-C2-O2
3	B	604	GOL	O1-C1-C2-O2
5	A	604	FBP	O5-C5-C6-O6
5	D	605	FBP	C4-C5-C6-O6
9	B	608	PEG	C4-C3-O2-C2

There are no ring outliers.

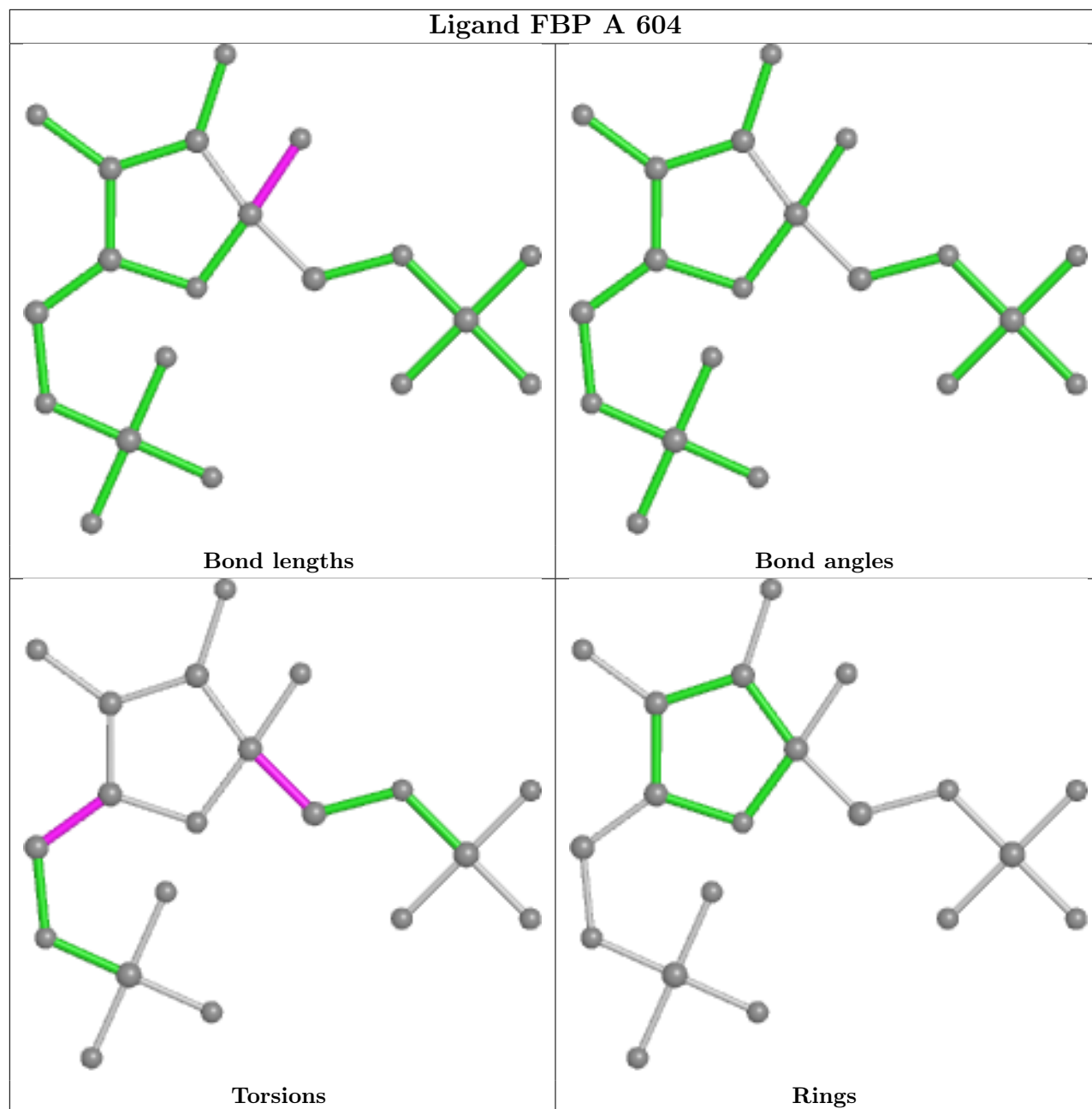
9 monomers are involved in 18 short contacts:

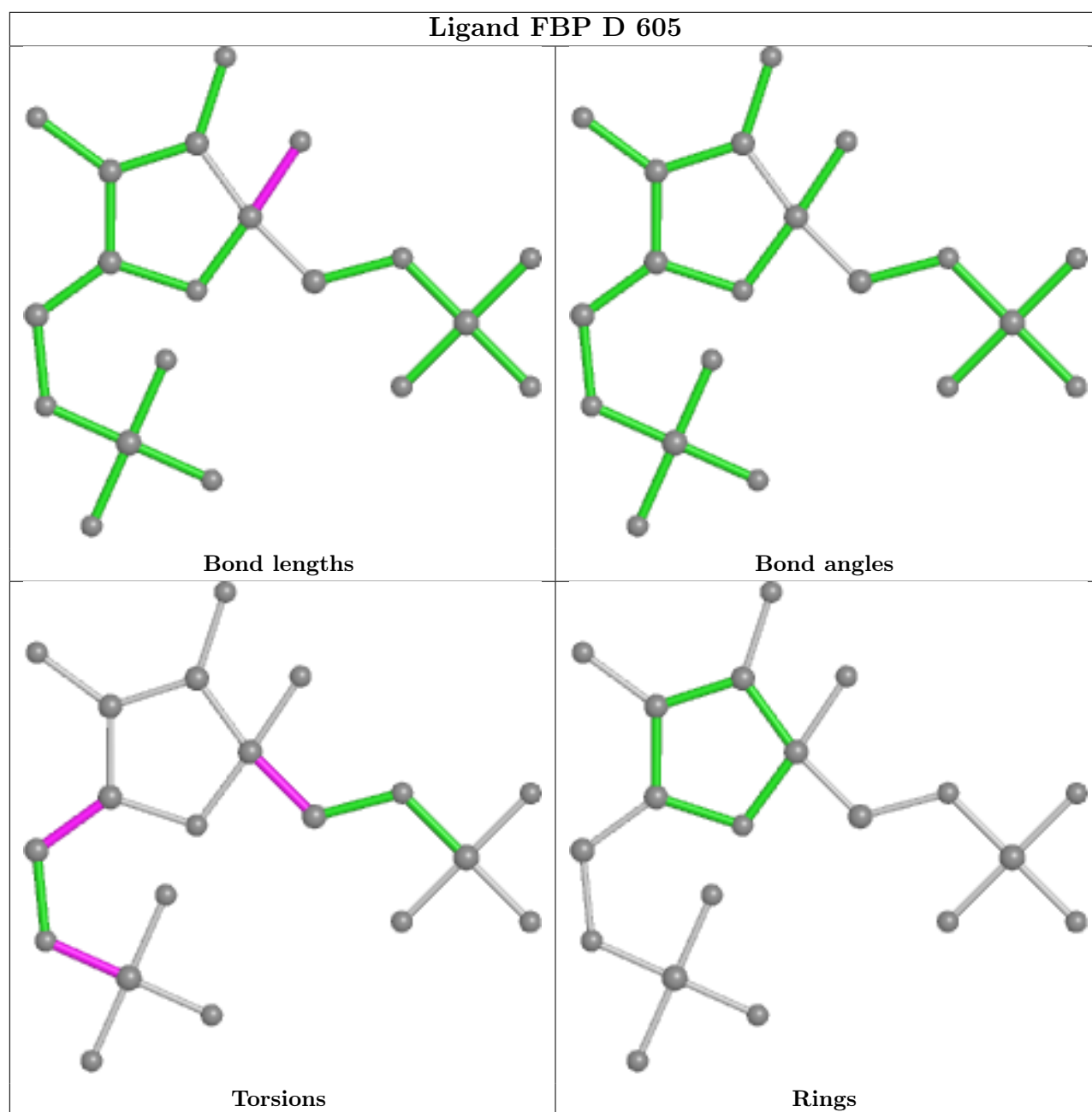
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	603	GOL	1	0
5	C	607	FBP	4	0
3	B	604	GOL	1	0
9	A	605	PEG	2	0
3	D	604	GOL	1	0
3	B	602	GOL	2	0
5	B	607	FBP	3	0
8	C	610	ASP	1	0
8	B	610	ASP	3	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.

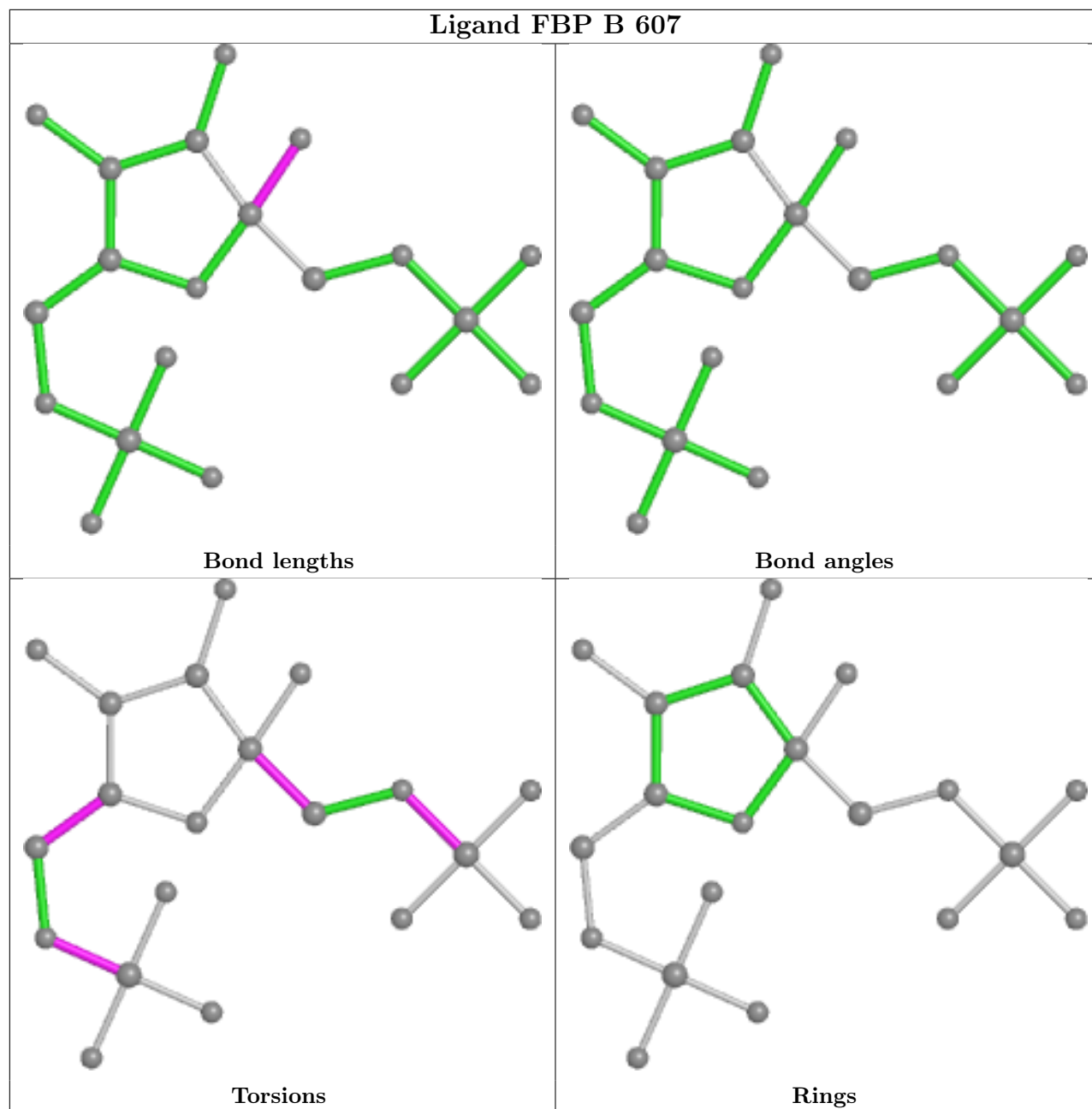


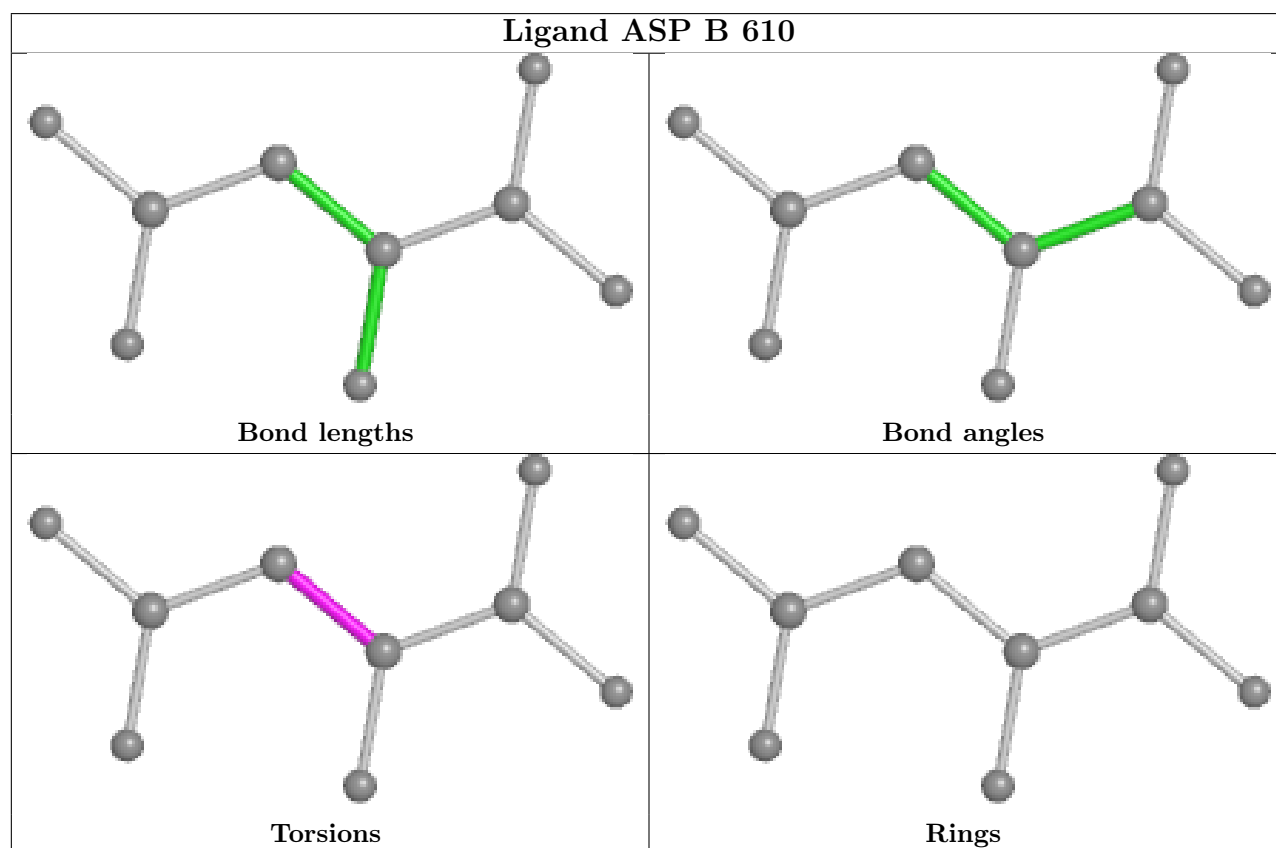
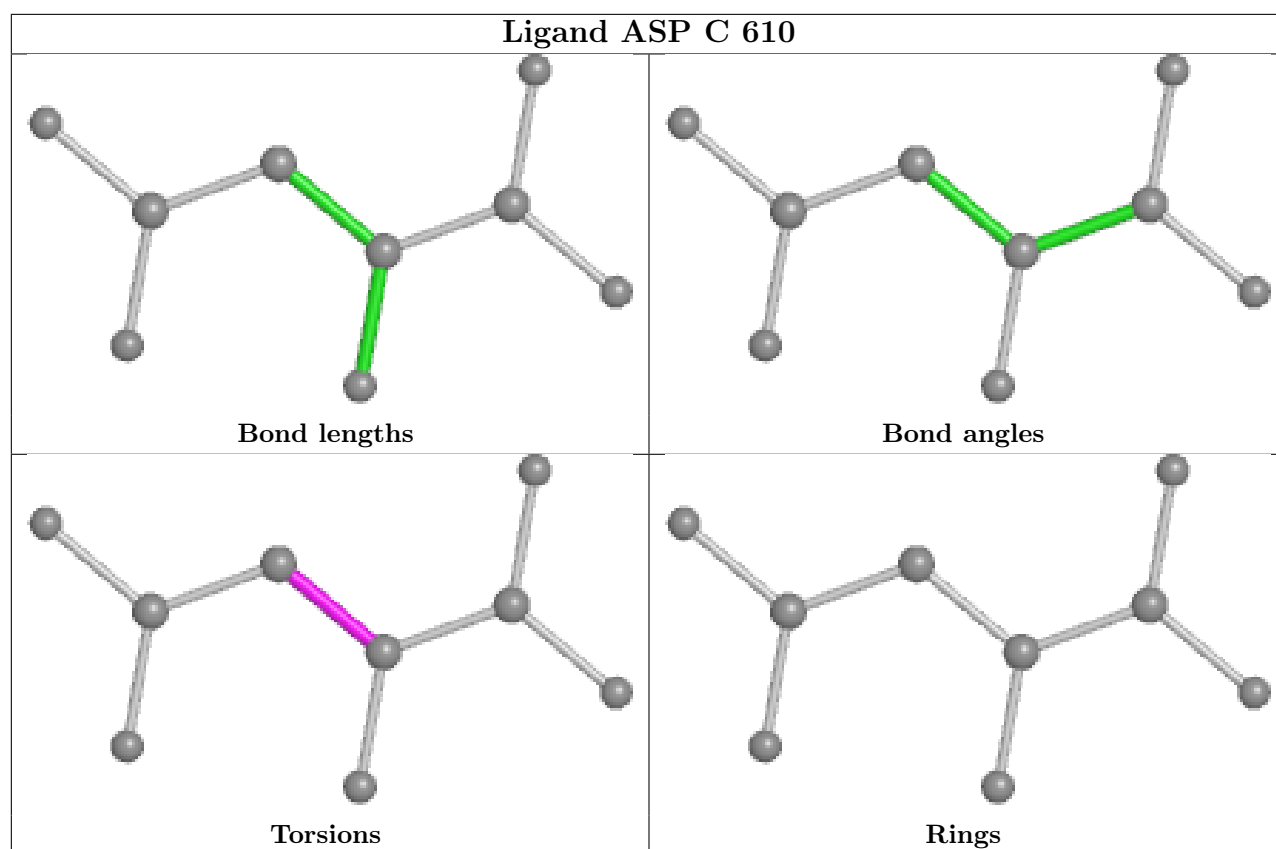
Ligand FBP A 604





Ligand FBP B 607





5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	509/550 (92%)	-0.23	4 (0%) 86 85	49, 67, 98, 114	0
1	B	512/550 (93%)	-0.11	9 (1%) 68 66	30, 66, 110, 129	0
1	C	517/550 (94%)	-0.15	8 (1%) 73 72	43, 59, 98, 115	0
1	D	430/550 (78%)	-0.26	3 (0%) 87 87	47, 64, 94, 131	0
All	All	1968/2200 (89%)	-0.18	24 (1%) 79 78	30, 64, 101, 131	0

All (24) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	163	ASN	3.4
1	A	100	SER	3.3
1	B	190	ALA	3.3
1	D	531	PRO	3.2
1	A	134	LEU	3.2
1	C	161	TYR	2.9
1	A	98	PHE	2.9
1	C	17	GLN	2.8
1	B	481	ALA	2.8
1	B	156	ILE	2.5
1	B	193	LEU	2.5
1	C	165	CYS	2.5
1	B	146	ASN	2.4
1	D	490	VAL	2.4
1	C	190	ALA	2.3
1	C	189	GLY	2.2
1	B	72	ALA	2.2
1	C	157	LEU	2.1
1	C	163	ASN	2.1
1	A	531	PRO	2.1
1	B	158	TRP	2.1

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Mol	Chain	Res	Type	RSRZ
1	D	216	VAL	2.1
1	C	166	LYS	2.1
1	B	187	GLN	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates [i](#)

There are no monosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
6	MG	A	607	1/1	0.71	0.42	60,60,60,60	0
9	PEG	B	608	7/7	0.72	0.28	70,74,79,86	0
6	MG	D	606	1/1	0.72	0.28	58,58,58,58	0
3	GOL	B	603	6/6	0.76	0.34	62,64,72,77	0
3	GOL	A	602	6/6	0.79	0.25	77,81,85,87	0
9	PEG	A	606	7/7	0.79	0.16	61,65,74,89	0
3	GOL	A	603	6/6	0.80	0.22	70,74,80,83	0
3	GOL	B	605	6/6	0.80	0.40	74,77,85,88	0
3	GOL	C	603	6/6	0.82	0.29	71,81,83,85	0
3	GOL	C	605	6/6	0.83	0.23	61,68,74,76	0
3	GOL	B	606	6/6	0.83	0.27	69,72,77,78	0
8	ASP	B	610	9/9	0.86	0.19	55,62,65,67	0
3	GOL	D	604	6/6	0.87	0.19	62,68,70,71	0
9	PEG	A	605	7/7	0.87	0.18	60,72,74,87	0
5	FBP	B	607	20/20	0.89	0.22	65,83,96,104	0
7	OXL	C	609	6/6	0.89	0.15	55,61,64,64	0
6	MG	B	609	1/1	0.89	0.30	59,59,59,59	0
2	K	C	602	1/1	0.89	0.08	82,82,82,82	0
3	GOL	C	604	6/6	0.91	0.40	50,59,60,76	0
2	K	D	602	1/1	0.91	0.17	114,114,114,114	0

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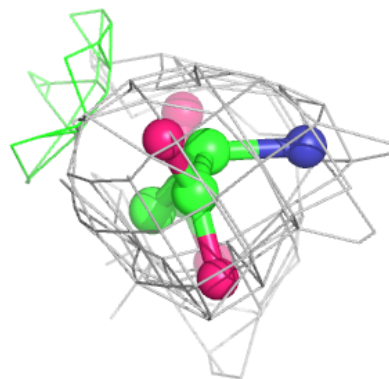
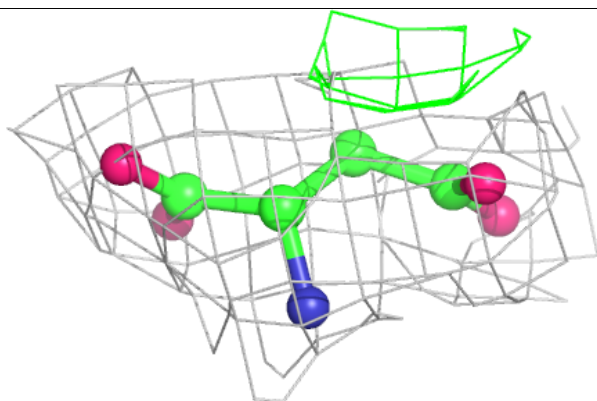
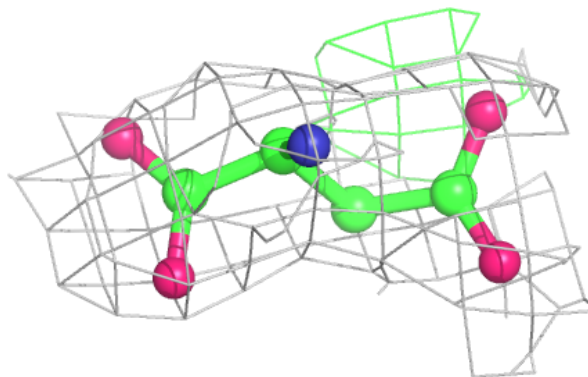
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	GOL	D	603	6/6	0.92	0.23	53,58,63,66	0
8	ASP	C	610	9/9	0.93	0.20	56,57,69,72	0
3	GOL	B	604	6/6	0.93	0.40	71,74,78,79	0
5	FBP	C	607	20/20	0.93	0.16	49,67,72,72	0
5	FBP	A	604	20/20	0.93	0.17	59,66,75,81	0
5	FBP	D	605	20/20	0.94	0.14	51,64,71,71	0
3	GOL	B	602	6/6	0.94	0.21	55,57,61,66	0
2	K	C	601	1/1	0.94	0.11	65,65,65,65	0
2	K	A	601	1/1	0.95	0.13	63,63,63,63	0
2	K	D	601	1/1	0.97	0.04	68,68,68,68	0
4	CL	C	606	1/1	0.98	0.13	47,47,47,47	0
6	MG	C	608	1/1	0.98	0.26	57,57,57,57	0
2	K	B	601	1/1	0.99	0.06	69,69,69,69	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

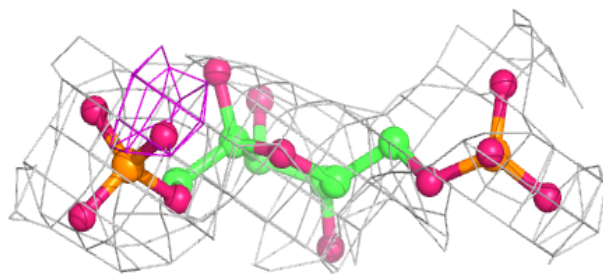
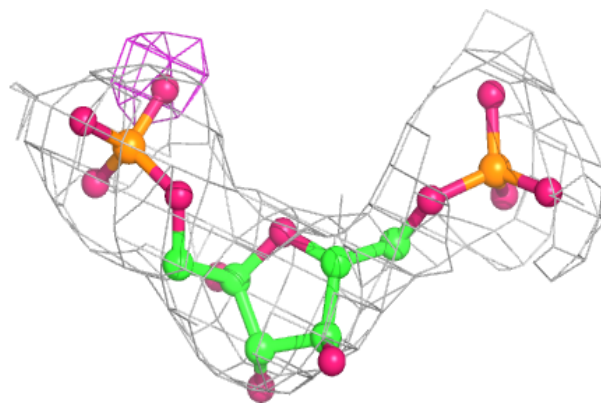
Electron density around ASP B 610:

2mF_o-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



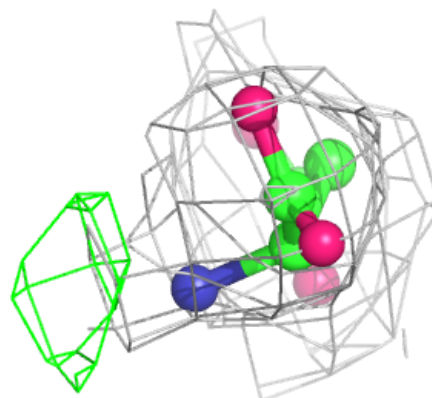
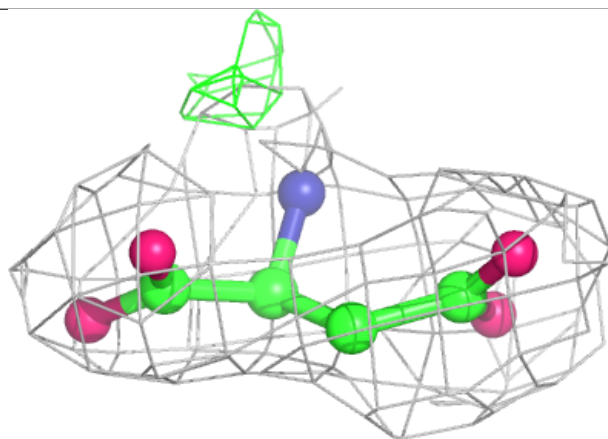
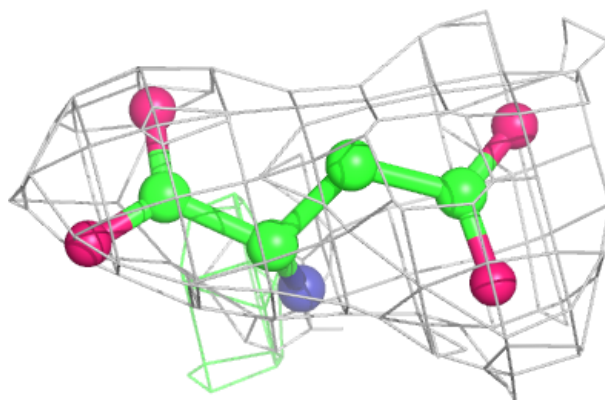
Electron density around FBP B 607:

$2mF_o - DF_c$ (at 0.7 rmsd) in gray
 $mF_o - DF_c$ (at 3 rmsd) in purple (negative)
and green (positive)



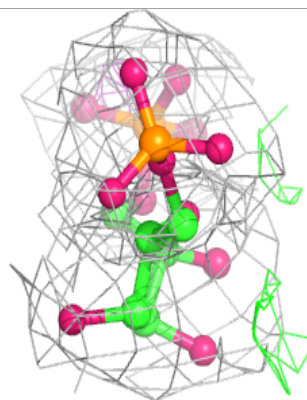
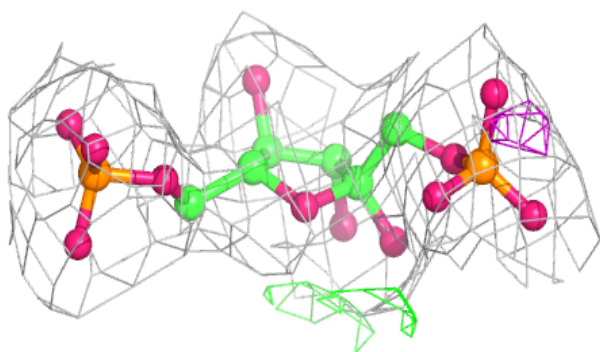
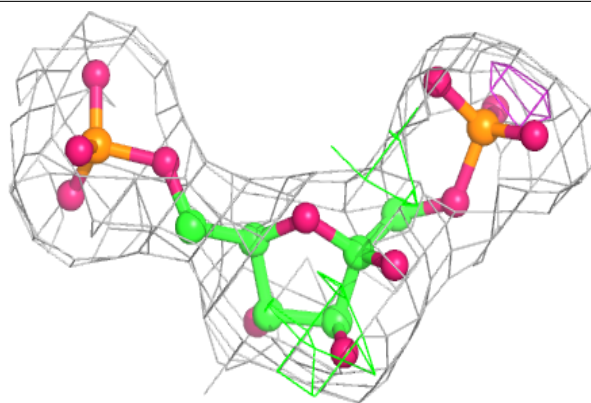
Electron density around ASP C 610:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

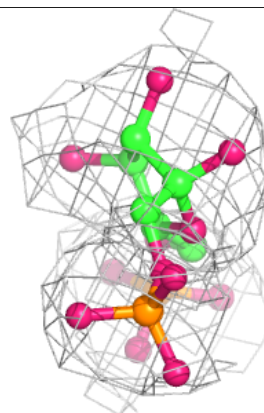
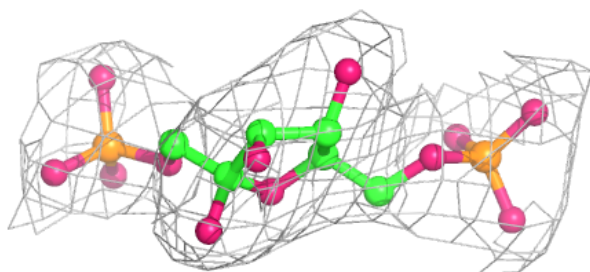
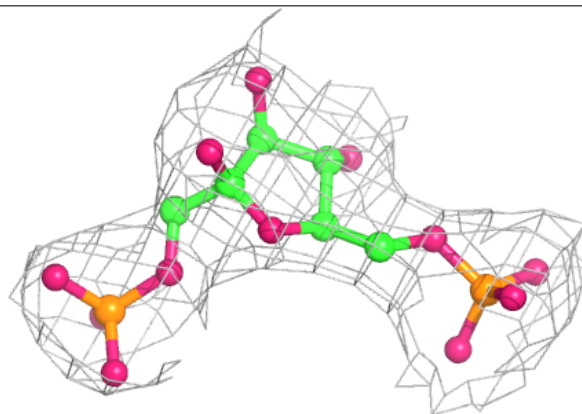


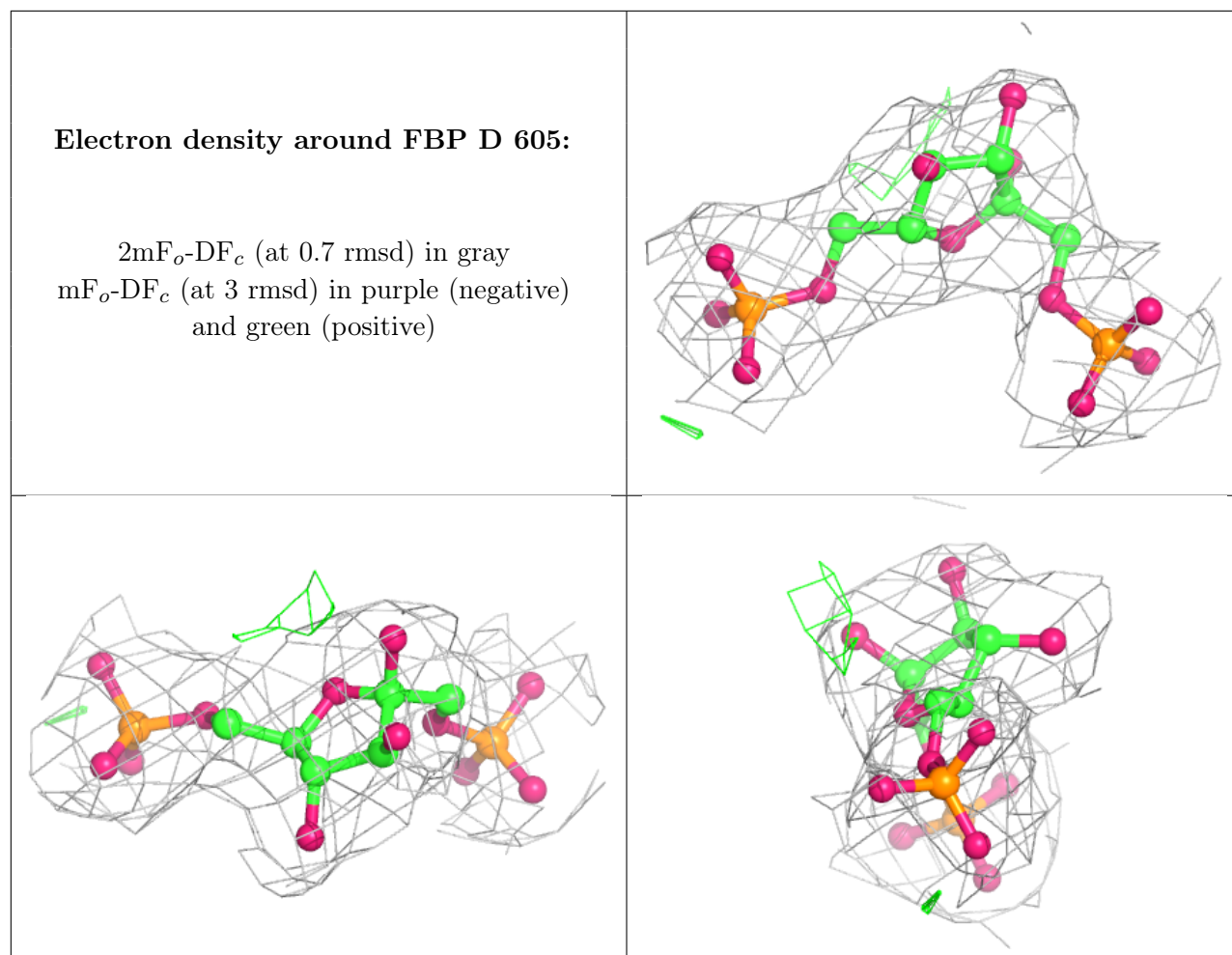
Electron density around FBP C 607:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

**Electron density around FBP A 604:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)





6.5 Other polymers [i](#)

There are no such residues in this entry.